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                 enhanced on STN
NEWS 4 JUN 26
                 NUTRACEUT and PHARMAML no longer updated
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         JUN 29
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NEWS 8 JUL 14 USGENE enhances coverage of patent sequence location
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NEWS 10 JUL 16 GBFULL adds patent backfile data to 1855
NEWS 11 JUL 21 USGENE adds bibliographic and sequence information
NEWS 12 JUL 28 EPFULL adds first-page images and applicant-cited
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NEWS 13 JUL 28 INPADOCDB and INPAFAMDB add Russian legal status data
NEWS 14 AUG 08 Improve STN by completing a survey and be entered to
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NEWS 16
         AUG 17 CAS REGISTRY, the Global Standard for Chemical
                 Research, Approaches 50 Millionth Registration
                 Milestone
NEWS 17 AUG 18 COMPENDEX indexing changed for the Corporate Source
                 (CS) field
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NEWS EXPRESS MAY 26 09 CURRENT WINDOWS VERSION IS V8.4, AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.

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=> FILE REG

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=>

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chain nodes : 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 31 32 33 ring nodes : 1 2 3 4 5 6 7 8 9 10 11 12 chain bonds : $1-18 \quad 2-25 \quad 3-19 \quad 4-26 \quad 5-13 \quad 6-24 \quad 7-21 \quad 8-14 \quad 9-22 \quad 10-23 \quad 11-17 \quad 12-20 \quad 13-14$ 13-16 14-15 17-27 18-32 19-31 27-29 27-28 28-33 ring bonds : 1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 8-9 9-10 10-11 11-12exact/norm bonds : 1-18 3-19 11-17 17-27 27-29 exact bonds : $2-25 \quad 4-26 \quad 5-13 \quad 6-24 \quad 7-21 \quad 8-14 \quad 9-22 \quad 10-23 \quad 12-20 \quad 13-14 \quad 13-16 \quad 14-15 \quad 18-32$ 19-31 27-28 28-33 normalized bonds : $1-2 \quad 1-6 \quad 2-3 \quad 3-4 \quad 4-5 \quad 5-6 \quad 7-8 \quad 7-12 \quad 8-9 \quad 9-10 \quad 10-11 \quad 11-12$

G1:Cy, Ak

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 31:CLASS 32:CLASS 33:CLASS

L1 STRUCTURE UPLOADED

=> D L1 L1 HAS NO ANSWERS L1 STR

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=> S L1 FULL FULL SEARCH INITIATED 16:19:54 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 447 TO ITERATE

100.0% PROCESSED 447 ITERATIONS 17 ANSWERS

SEARCH TIME: 00.00.01

L2 17 SEA SSS FUL L1

=> FILE CAPLUS

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 186.36 186.58

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FILE COVERS 1907 - 20 Aug 2009 VOL 151 ISS 8

FILE LAST UPDATED: 19 Aug 2009 (20090819/ED)

REVISED CLASS FIELDS (/NCL) LAST RELOADED: Jun 2009

USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Jun 2009

CAplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

The ALL, BIB, MAX, and STD display formats in the CA/CAplus family of databases have been updated to include new citing references information. This enhancement may impact record import into database management software. For additional information, refer to NEWS 9.

=> S L2 L3 2 L2

=> D L2 IBIB ABS HITSTR 1-2
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'ABS' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'
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The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

EPROP - Table of experimental properties PPROP - Table of predicted properties PROP - EPROP, ETAG, PPROP and SPEC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original) OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

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=> D L3 IBIB ABS HITSTR 1-2

ANSWER 1 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2008:1530346 CAPLUS

DOCUMENT NUMBER: 150:77699

TITLE: Compositions and methods of use for treating or

preventing lipid related disorders

INVENTOR(S): Currie, Mark; Talley, John; Cali, Brian

PATENT ASSIGNEE(S): Ironwood Pharmaceuticals, Inc, USA

SOURCE: PCT Int. Appl., 197pp. CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PA	PATENT NO.				KIND		DATE		APPLICATION NO.									
	2008	2008157537 2008157537					20081224		WO 2008-US67204					20080617				
WO							2009											
	W:	ΑE,	AG,	AL,	ΑM,	ΑO,	ΑT,	ΑU,	AZ,	BA,	BB,	BG,	BH,	BR,	BW,	BY,	BZ,	
		CA,	CH,	CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DO,	DZ,	EC,	EE,	EG,	ES,	
		FΙ,	GB,	GD,	GE,	GH,	GM,	GT,	HN,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	ΚE,	
		KG,	KM,	KN,	KP,	KR,	KΖ,	LA,	LC,	LK,	LR,	LS,	LT,	LU,	LY,	MA,	MD,	
		ME,	MG,	MK,	MN,	MW,	MX,	MY,	MΖ,	NA,	NG,	NI,	NO,	NZ,	OM,	PG,	PH,	
		PL,	PT,	RO,	RS,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SM,	SV,	SY,	TJ,	TM,	
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		IE,	IS,	IT,	LT,	LU,	LV,	MC,	MT,	NL,	NO,	PL,	PT,	RO,	SE,	SI,	SK,	
											GQ,		•				TD,	
		TG,	BW,	GH,	GM,	KE,	LS,	MW.	ΜZ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	
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OTHER SOURCE(S): MARPAT 150:77699

GΙ

Disclosed herein are compds. of formula I and II and their compns. and AΒ methods for treating or preventing a variety of disorders and conditions associated with lipid metabolism The methods generally include administering to

a patient in need thereof a therapeutically effective amount of a pharmaceutical composition comprising one or more fibric acid or statin derivative

compns. alone or in combination with one or more lipid altering agents and/or PDE inhibitors. Compds. of formula I and II wherein R1 is H and halo; R2 is H, halo, (un) substituted cycloalkyl, (un) substituted benzoyl, etc.; Z is O, and (CH2)1-3-O; X is a bond, O, NH, and amino acid residue; R4 is Oh, NO, NO2, amino acid residue, fibric acid residue, guanidine, tetrazolyl, agmatine, etc.; R5 is a statin residue; are claimed. Example compound III was prepared by a general procedure. The invention compds. were evaluated for their ability to treat lipid related disorders.

ΙT 1094098-94-8P

> RL: PAC (Pharmacological activity); PKT (Pharmacokinetics); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of compds. for treatment, prevention and combination therapy of lipid-related disorders)

RN 1094098-94-8 CAPLUS

Propanoic acid, 2-[4-(4-chlorobenzoyl)phenoxy]-2-methyl-, CN 4-[(1E)-2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

Double bond geometry as shown.

L3 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2005:696641 CAPLUS

DOCUMENT NUMBER: 143:172689

TITLE: Preparation of resveratrol ester analogs as sirtuin

activators

INVENTOR(S): Andrus, Merritt B.; Liu, Jing

PATENT ASSIGNEE(S): Brigham Young University Technology Transfer Office,

USA

SOURCE: PCT Int. Appl., 95 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.				KIND		DATE			APPLICATION NO.				DATE					
WO 2005069998				A2		20050804			WO 2005-US2229					20050119				
WO	WO 2005069998			A3 20060105			0105											
	W:	ΑE,	AG,	AL,	AM,	ΑT,	ΑU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,	
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,	
		LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MΖ,	NA,	ΝΙ,	
		NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SY,	
		TJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC,	VN,	YU,	ZA,	ZM,	ZW,	SM
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AU 2005207029			•					AU 2005-207029				20050119						
			A1 20060804				CA 2005-2593576					20050119						
EP 1753708							EP 2005-711939											
										EE,								
	•						,			PT,					TR	/	,	

US 20080255382 A1 20081016 US 2006-597335 20060721
PRIORITY APPLN. INFO.: US 2004-537622P P 20040120
US 2004-616537P P 20041006
WO 2005-US2229 W 20050119

OTHER SOURCE(S): CASREACT 143:172689; MARPAT 143:172689

Т

GΙ

AB Resveratrol and ester analogs of formula I [A1-A3 = protecting group, acyl] are prepared The compds. are made from a multi-step process including a N-heterocyclic carbene-type ligand coupling in the presence of a base with benzoyl halide and styrene coupling partners. These compds. show increased stability for use in the food, cosmetic and pharmaceutical industries (no data). Thus, resveratrol (I; A1-A3 = H) was prepared by decarbonylative Heck coupling of 3,5-diacetoxybenzoyl chloride using Pd(OAc)2 and 1,3-bis(2,6-diisopropylphenyl)imidazolinium chloride and 3-acetoxystyrene followed by deprotection with NaOH.

861446-31-3P 861446-36-8P 861446-41-5P ΤТ 861446-46-0P 861446-51-7P 861446-56-2P 861446-61-9P 861446-66-4P 861446-71-1P 861446-76-6P 861446-81-3P 861446-86-8P 861446-96-0P 861447-01-0P 861446-91-5P 861447-06-5P

RL: COS (Cosmetic use); FFD (Food or feed use); PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of resveratrol ester analogs as sirtuin activators)

RN 861446-31-3 CAPLUS

CN 1,3-Benzenediol, 5-[2-[4-(1-oxopropoxy)phenyl]ethenyl]- (CA INDEX NAME)

RN 861446-36-8 CAPLUS

CN Butanoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

RN 861446-41-5 CAPLUS

CN Pentanoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

RN 861446-46-0 CAPLUS

CN Hexanoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

RN 861446-51-7 CAPLUS

CN 2,4-Hexadienoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

RN 861446-56-2 CAPLUS

CN Dodecanoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

HO CH CH CH O CH O CH
$$_{0}$$
 CH $_{0}$ CH $_{0}$ Me

RN 861446-61-9 CAPLUS

CN Hexadecanoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

RN 861446-66-4 CAPLUS

CN Octadecanoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

RN 861446-71-1 CAPLUS

CN 9-Octadecenoic acid (9Z)-, 4-[(1E)-2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

Double bond geometry as shown.

Me
$$(CH_2)$$
 7 Z (CH_2) 7 O

RN 861446-76-6 CAPLUS

CN 9,12-Octadecadienoic acid (9Z,12Z)-, 4-[(1E)-2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

Double bond geometry as shown.

PAGE 1-A OH (CH2)
$$\frac{1}{4}$$
 $\frac{1}{2}$ $\frac{1}{2}$ (CH2) $\frac{1}{2}$ O

PAGE 1-B

OH

RN 861446-81-3 CAPLUS

CN 6,9,12-Octadecatrienoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

- CH $_2$ - CH= CH- (CH $_2$) $_4$ - Me

RN 861446-86-8 CAPLUS

CN 9,12,15-Octadecatrienoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

- CH $_2$ - CH= CH- Et

RN 861446-91-5 CAPLUS

CN 3,6,9-Octadecatrienoic acid, 4-[2-(3,5-dihydroxyphenyl)] ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

-CH=CH-(CH₂) $_{7}-$ Me

RN 861446-96-0 CAPLUS

CN 5,8,11,14-Eicosatetraenoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

$$-$$
 CH $_2-$ CH $-$ CH $_2-$ CH $-$ CH $_2-$ CH $-$ (CH $_2$) $_4-$ Me

RN 861447-01-0 CAPLUS

CN 5,8,11,14,17-Eicosapentaenoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

$$-$$
 CH₂ $-$ CH $-$ Et

RN 861447-06-5 CAPLUS

CN 4,7,10,13,16,19-Docosahexaenoic acid, 4-[2-(3,5-dihydroxyphenyl)ethenyl]phenyl ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

OS.CITING REF COUNT: 6 THERE ARE 6 CAPLUS RECORDS THAT CITE THIS RECORD

(6 CITINGS)

REFERENCE COUNT: 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS

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